# MEMORANDUM (LABORATORY DATA REPORT)

**EPA - General Parameters** 

In reply refer to: 10-LC94

To: Rick Wilkin From: Lynda Callaway

Lab: General Parameters

Thru: Cindy Paul Date: 10/20/2010

Mark White Kristie Hargrove

Work Request: EPAGP216 Copies: Rick Wilkin

Task No.: 23993 Cindy Paul

Kristie Hargrove Lynda Callaway Steve Vandegrift

Sample Site/Project: Pavillion Groundwater

Date Received: 10/8/2010 Sample Matrix: Groundwater

Date Analyzed: 10/13/2010 Analysis Type: Chloride, sulfate & fluoride

No. Samples Analyzed: 9 Sample Preparation: Diluted as needed

Method(s) Used: RSKSOP-276, Rev. 3 - Determination of Major Anions in Aqueous Samples Using Capillary

Electrophoresis With Indirect UV Detection and Empower 2 Software

## Comments:

Quality control measures performed along with your samples included analysis of method blanks, sample matrix spikes, laboratory sample duplicates, calibration check standards, and second-source quality control samples as outlined in RSKSOP-276, revision 3. Method detection limits (MDLs) were determined on 10/1/2010. Note that a sample for anions was not received for field sample ID LD 02 dup.

# EPA - General Parameters Analytical Results Report

Laboratory:	General Parameters	]							
Work Request:	EPAGP216	Sample Results							
Analyst: Lynda Callaway		]	Analytes	Chloride (Cl <sup>-</sup> )		Sulfate ( SO <sub>4</sub> <sup>-2</sup> )		Fluoride (F <sup>-</sup> )	
			Codes	16887-00-6		14808-79-8		7782-41-4	
Report Date:	10/20/10	]	Methods	RSKSOP-276/3		RSKSOP-276/3		RSKSOP-276/3	
		_	Unit	mg/L		mg/L		mg/L	
			MDL	* 0.136		* 0.103		* 0.056	
			QL	* 1.00		* 1.00		* 0.200	
Field Sample ID	Lab Sample ID	Date Collected	Date Analyzed	Data	DF	Data	DF	Data	DF
RD 01	5763-1	10/5/2010	10/13/2010	15.2	1	357	21	0.992	1
EPAMW 01	5763-2	10/6/2010	10/13/2010	23.3	1	398	21	1.55	1
EPAMW 02	5763-3	10/6/2010	10/13/2010	466	21	12.1	1	1.01	1
LD 01	5763-4	10/6/2010	10/13/2010	33.0	4	1,320	50	0.898	4
LD 01	5763-4 Lab dup	10/6/2010	10/13/2010	32.9 (RPD=0.303)	4	1,310 (RPD=0.760)	50	0.926 (RPD=3.07)	4
LD 01 (Dup)	5763-5	10/6/2010	10/13/2010	32.9	4	1,320	50	0.988	4
RD 01 Field Blank	5763-6	10/5/2010	10/13/2010	ND	1	ND	1	ND	1
Trip Blank	5763-7	10/6/2010	10/13/2010	ND	1	ND	1	ND	1
EQ Blk	5763-8	10/7/2010	10/13/2010	ND	1	ND	1	ND	1
LD 02	5763-9	10/7/2010	10/13/2010	20.1	1	698	100	2.28	2
LD 02	5763-9 Lab dup	10/7/2010	10/13/2010	19.7 (RPD=2.01)	1	691 (RPD=1.01)	100	2.17 (RPD=4.94)	2
LD 02 dup	5763-10	10/7/2010	1	-	-	-	-	-	-

### Comments:

The data quality objective for the precision of sample duplicates is a relative percent difference (RPD) of < 10%. This objective was met for all samples within the range of the calibration standards. MDLs were determined on 10/1/2010. \* The MDLs and QLs should be raised by the same factor as the dilution factor for those samples that were diluted. Note that an anion sample was not received for field sample ID LD 02 dup.

### Notes

<sup>1.</sup> If the parameter was detected above the quantitation limit (QL), the numeric result is reported; BQL denotes that the parameter was not detected at or above the quantitation limit; BQL () denotes that the parameter was detected above the method detection limit (MDL) but below QL and the estimated numeric result is reported in parenthesis; ND denotes that the parameter was not detected at all. All the results are corrected with dilution factors (DF), if applicable. NA means not applicable.

<sup>2. &</sup>quot;-" denotes that the information is not available or the analyte is not analyzed.

# **EPA - General Parameters** Analytical Results Report

General Parameters Laboratory:

EPAGP216 Work Request:

Quality Control Data Summary

Analyst:	Lynda Callaway					
Report Date:	10/20/10					

Additional ID

RO water Blank

ERA #46 Minerals

Calibration Check Standard

EPAMW 01 Spike

Date Prepar

QC Sample ID

MB

MB

MB

MB

MB

MB

SS

SS

SS

SS

SS

ccc

CCC

CCC

CCC

CCC

CCC

CCC

ccc

CCC

MS

				2							
	Analytes	Chloride (Cl')			Sulfate (SO <sub>4</sub> <sup>-2</sup> )			Fluoride (F')			
	Codes         16887-00-6           Methods         RSKSOP-276/3			14808-79-8 RSKSOP-276/3			7782-41-4				
							RSKSOP-276/3				
	Unit	Unit mg/L MDL 0.136			mg/L 0.103			mg/L 0.056			
	MDL										
	QL	1.00			1.00			0.200			
Date Prepared	Date Analyzed	Data	True Value	% REC.	Data	True Value	% REC.	Data	True Value	% REC.	
10/13/2010	10/13/2010	ND	-		ND	-	-	ND		-	
10/13/2010	10/13/2010	ND	-		ND	-	-	ND	-	-	
10/13/2010	10/13/2010	ND	-	-	ND	-	-	ND	-	-	
10/13/2010	10/13/2010	ND	-	-	ND	-	-	ND	-	-	
10/13/2010	10/13/2010	ND	-	-	ND	-	-	ND	-	-	
10/13/2010	10/13/2010	ND	-		ND	-		ND		-	
6/24/2010	10/13/2010	51.3	52.0	98.7	26.8	28.5	94.0	2.69	2.70	99.6	
6/24/2010	10/13/2010	52.6	52.0	101	27.6	28.5	96.8	2.74	2.70	101	
6/24/2010	10/13/2010	53.0	52.0	102	28.0	28.5	98.2	2.69	2.70	99.6	
6/24/2010	10/13/2010	52.7	52.0	101	27.6	28.5	96.8	2.75	2.70	102	
6/24/2010	10/13/2010	53.1	52.0	102	28.1	28.5	98.6	2.69	2.70	99.6	
10/1/2010	10/13/2010	1.05	1.00	105	BQL (0.916)	1.00	91.6	0.210	0.200	105	
10/1/2010	10/13/2010	5.14	5.00	103	5.05	5.00	101	1.04	1.00	104	
10/1/2010	10/13/2010	25.3	25.0	101	25.2	25.0	101	5.07	5.00	101	
10/1/2010	10/13/2010	1.02	1.00	102	BQL (0.935)	1.00	93.5	0.205	0.200	103	
10/1/2010	10/13/2010	5.13	5.00	103	5.09	5.00	102	0.963	1.00	96.3	
10/1/2010	10/13/2010	25.8	25.0	103	25.9	25.0	104	5.19	5.00	104	
10/1/2010	10/13/2010	5.22	5.00	104	5.13	5.00	103	1.06	1.00	106	
10/1/2010	10/13/2010	5.19	5.00	104	4.98	5.00	99.6	1.00	1.00	100	
10/1/2010	10/13/2010	25.4	25.0	102	25.4	25.0	102	5.03	5.00	101	
10/13/2010	10/13/2010	32.7	23.3 (9.62)	97.7	* 27.7	* 19.0 (9.62)	90.4	3.53	1.55 (1.92)	103	

The data quality objective (DQO) for the accuracy of continuing calibration check standards is 90-110% recovery. The DQO for ERA # 46 is 84.8 - 116% for chloride, 79.6 - 118% for sulfate, and 83.0 - 117% for fluoride. The DQO for the recovery of matrix spikes is 80 - 120%. These objectives were met for the standards and spikes. The matrix spikes were prepared by adding 20 uL of a 250 / 50 mg/L mixed standard into 0.5 mL of sample yielding a spike concentration of 9.62 mg/L for chloride and sulfate and 1.92 mg/L for fluoride. The matrix spike recovery was calculated according to the equation: % Recovery = 100 x (Spiked sample concentration (DATA) - Native sample concentration //Spike concentration. \*The values for spike concentrations are calculated and reported without the dilution factors applied.

1. MB - Method Blank, CCC - Continuing Calibration Check. A calibration standard analyzed within the batch of samples, LCS - Laboratory Control Spike. A laboratory blank spiked with analytes at known concentrations. MS - Matrix Spike. A field sample spiked with known concentrations of analytes. The field sample id is identified. The True Value column for matrix spikes list the unspiked native sample concentration along with the spike concentration in parentheses. SS - Samples obtained from the second sources are identified by their designated names. DUP - Field sample duplicate analysis. A sample selected by the lab analyst to analyze as a duplicate. It is reported in the sample result section. % REC - Percent Recovery. Calculated as the percentage of the results to the true values. It equals to % accuracy for CCC.